A cross-institutional, FAIR VIVO for Metabolomics

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Metabolomics

Metabolomics is the scientific study of metabolites present within an organism, cell, or tissue. Human metabolites are small molecules found in human tissue that occur naturally as a result of human metabolism, or are present as a result of drugs, food components, or exposure to environmental conditions. Along with genomics (the study of DNA), transcriptomics (RNA), proteomics (proteins), metabolomics provides information regarding biochemical compounds and processes in cells, leading to a better understanding of cellular biology. "Metabolic profiling can give an instantaneous snapshot of the physiology of a cell, and thus, metabolomics provides a direct "functional readout of the physiological state" of an organism." (Wikipedia http://bit.ly/2Pg06X3)

Such profiles can be used to distinguish tissue types, disease states, and health status of individuals. Understanding the changes in profiles over time can lead to improved understanding of diseases such as cancer and diabetes, leading to potential treatments.

Over 450,000 human metabolites have been identified to date. Metabolomics is considered an

Metabolomics Data

The Metabolomics Workbench (MWB) (https://metabolomicsworkbench.org) is the National Metabolomics Data Repository of the US National Institutes of Health (NIH). Anyone can deposit data to the MWB. As of August 28, 2019, the workbench provides data from 982 publicly available studies. Another 205 studies are currently embargoed and will be available subject to their embargo dates. MWB develops and uses RefMet (http://bit.ly/2PkxY5p), a nomenclature for representing metabolites found using mass spectroscopy (MS) and nuclear magnetic resonance (NMR) techniques. Investigators upload study data to the workbench and provide metadata regarding themselves and their work. The MWB provides an API that can be used to access metadata about studies and investigators. MWB metadata has been mapped to an ontology developed by the authors to represent it as RDF and load it to VIVO.

PubMed (<u>http://pubmed.gov</u>) is an open access index to literature in metabolomics. Its API can be used to find and retrieve publication data regarding metabolomics investigators. Many groups use PubMed data in VIVO. PubMed data has been mapped to the VIVO Ontology (<u>http://vivoweb.org/ontology/core</u>)

FAIR Data Principles

The FAIR Data Principles (Wilkinson, et. al <u>https://www.nature.com/articles/sdata201618</u>) provide a framework for creating data that can be used by groups beyond the group that created the data. The principles are difficult to achieve in practice and much has been written about implementation. VIVO supports the principles naturally, it is designed to share data.

- Findable data should be found using search engines on the Internet
- Accessible the data can be accessed without technical, legal or operational barriers
- Interoperable the data is available in a common format, so that data from more than one source can be readily combined
- Reusable the data can be reused by those finding it. This typically means that enough information is provided for the recipients of the data to determine if thre data is suitable for their purposes.

The Metabolomics Workbench (<u>https://metabolomicsworkbench.org</u>), Metabolites (<u>https://www.ebi.ac.uk/metabolights/index</u>), and the MassBank (<u>http://massbank.jp</u>) provide data

emerging field. Technical challenges in compound identification and data analysis are significant.

The NIH Common Fund Metabolomics consortium originated in 2013 to help address these and other challenges. Fourteen investigators across the US are supported to advance metabolomics. Five are engaged in compound identification, seven in software tools and data analysis. The group at UCSD supports the Metabolomics Workbench, a data repository. The group at the University of Florida provides overall coordination for the consortium.

Data regarding software used in metabolomics studies has been difficult to find. An index to such software will be created and curated by a group at the University of Colorado Anschutz Medical Campus funded by the NIH. This data will be in the form of a spreadsheet. The authors will use the Software Ontology (SWO) (http://www.obofoundry.org/ontology/swo.html) to represent the data as RDF and load the data into VIVO.

on experimental results, and spectra of compounds according to the FAIR principles. In each case, data can be found by using internal search capabilities of each of the sites. This is not ideal. Generally findable results using search engines can be created for metabolomic data sets using the techniques described here. Each dataset or study will have its own page which can be found and indexed by search engines. Our approach will result in pages for datasets, publications, investigators, and software.

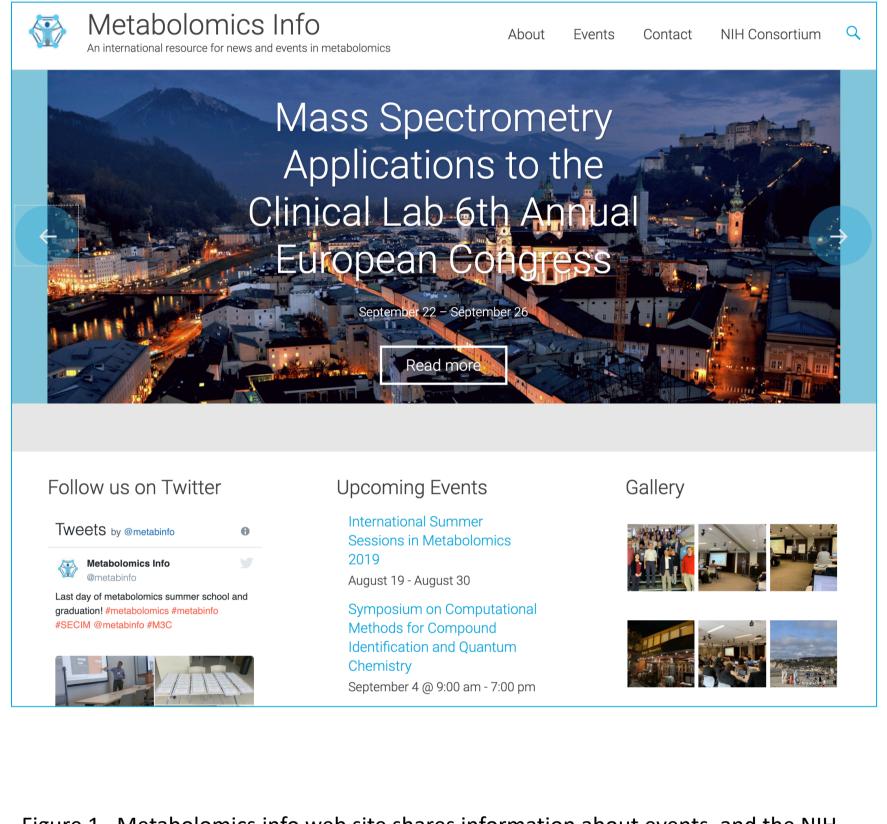


Figure 1.	Metabolomics.info	veb site shares	information	about events,	and the NIH
consortiu	m				

METABOL	OMICS	Log in / Registe
WORKBE	NCH	Search the Metabolomics Workbench
ne NIH Data Repository Databases Protocols Standard		
NIH Metabolomics Data Repository	Quick Links - Key Resources	
Upload and Manage Browse and Search Studies Studies	Analyze Studies	EVENTS CALENDAR
As of 08/28/2019 the NIH Metabolomics Data Repository contains 982 public A total of 1187 studies have been processed by the DRCC and the remainder subject to their embargo dates.	Regional Comprehensive Metabolomics Resource Cores (RCMRC)s	
		Michigan Regional Comprehensive Metabolomics Resource Core (MRC) ² 립
Metabolomics News	NIH West Coast Metabolomics Center at UC Davis 妤	
08-02-2019: RefMet: A <u>Ref</u> erence list of <u>Met</u> abolite names	Eastern Regional Comprehensive Metabolomics Resource Core (ERCMRC) &	
The main objective of RefMet E is to provide a standardized reference nom metabolite structures and isobaric species identified by MS and NMR techniq	Southeast Center for Integrated Metabolomics (SECIM) &	
This is an essential prerequisite for the ability to compare and contrast experiments and studies. To this end, a list of over 200,000 names from a s	Resource Center for Stable Isotope- Resolved Metabolomics (RC-SIRM)	
studies on the Metabolomics Workbench has been used as a starting point to analytical chemistry-centric list of common names for metabolite structures	Metabolomics Core at Mayo Clinic 🗗	
these names have been linked to a metabolite classification system using and ClassyFire regional classification methods. A name-conversion user interface	▹ more about our RCMRCs	
a list of metabolite names and map them to the corresponding Refmet names increase the data-sharing potential of metabolomics experiments and fac biology objectives for the majority of commonly encountered metabolite spec	YOUR FEEDBACK- Please tell us about your experiences with the Metabolomics Workbench.	
 Browse/Search/Download Refmet Convert metabolite names to RefMet nomenclature 		
 Help on RefMet 		
▶ news archive		
UCSD Metabolomics Workbench, a resource sponsored by the Common Fund		map Contact Personnel (Restricted access)
UC San Diego		

Figure 2. The Metabolomics Workbench is a data repository for data from metabolomics experiments. An API provides metadata regarding studies and investigators for all to use.

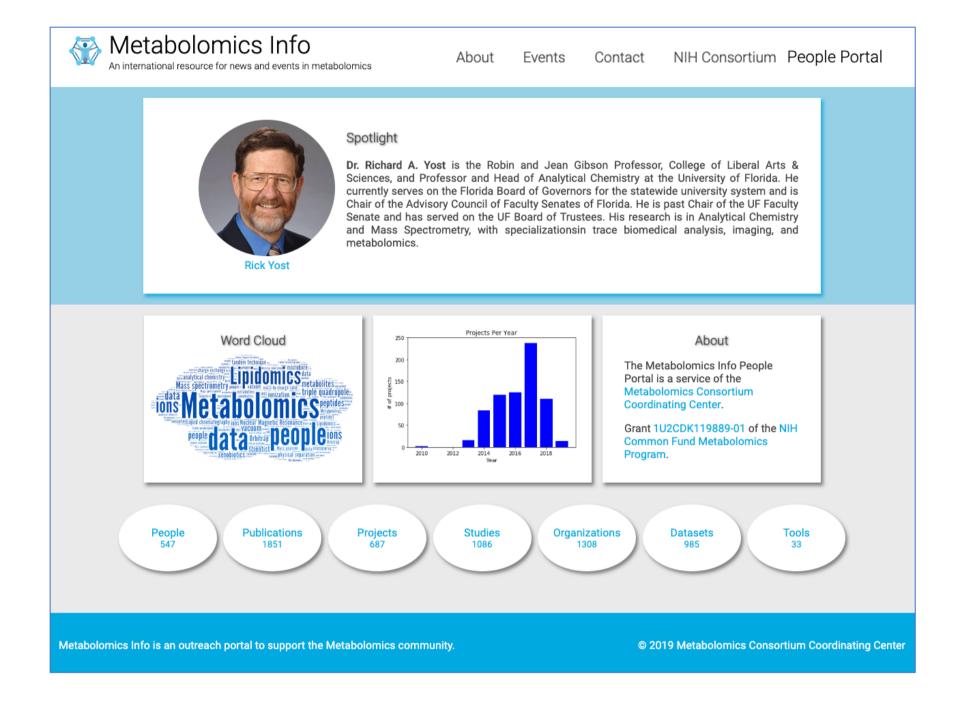


Figure 1. Metabolomics.info web site shares information about events, and the NIH consortium

Use Cases

The metabolomics community is still developing. Some use cases for the web site being developed include:

- 1. Discover metabolomics datasets using search engines. The pages created by the approach described here.
- Organize metabolomics information by investigator. Creating web pages for each investigator will emphasize the number of datasets deposited by each investigator, the number of metabolomics papers of the investigator, and software systems created by the investigator.
- 3. Provide summaries of contributions by investigator.
- 4. Provide contact information for investigators.
- 5. Boost the visibility of the work products (papers, software, and datasets) of the NIH Common Fund Metabolomics Consortium, leading to increased collaboration between consortium members and others around the world.
- 6. Increase deposits to Metabolomics workbench
- 7. Increase reuse of data in Metabolomics workbench
- 8. Through the FAIR data principles, encourage data sharing and reuse in the metabolomics community.

Triple Pattern Fragments (TPF)

Triple Pattern Fragments (TPF), (Verborgh, et. Al doi:10.1016/j.websem.2016.03.003) is a low cost knowledge interface for the web. TPF provides a simple, fast, general mechanism for querying knowledge graphs of RDF triples. TPF answers queries of the form:

Subject predicate object

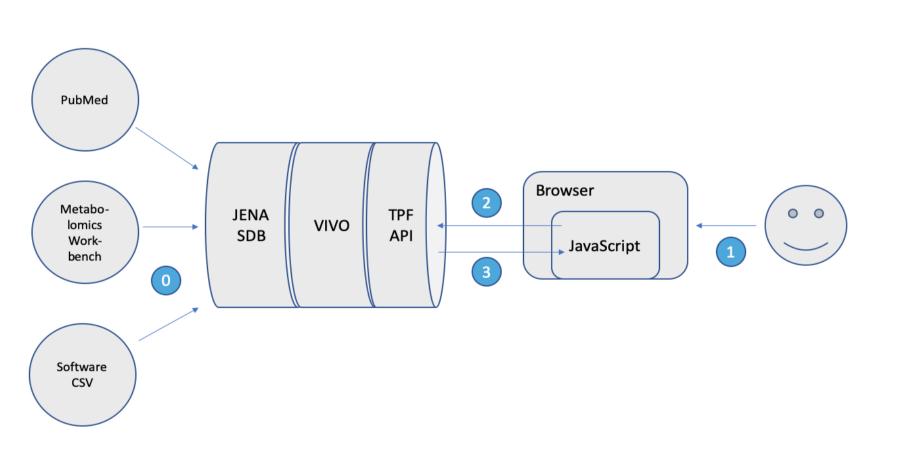
Where any of the three components can be a wildcard. A query such as

* rdf:type *

returns all the type assertions in a triple store. When executed at http://openvivo.org/tpf/core,

Technical Approach

See figure below. 0) Publication data from PubMed, dataset metadata from Metabolomics workbench, and software metadata from collaborators at the University of Colorado Anschutz Medical Campus are retrieved, transformed to triples and loaded to the JENA SDB triple store (part of VIVO) using custom scripts on a regular basis; 1) User makes request for page that will contain data from VIVO; 2) TPF Client JavaScript software makes a TPF query of the VIVO TPF endpoint. The JavaScript software is generic – any ontology and any TPF endpoint can be used; 3) the TPF API returns a set of triples to the client. No modifications of VIVO are necessary.



JavaScript

A small JavaScript library is available (<u>http://github.com/ctsit/tpf</u>) to make TPF queries and handle TPF replies. The code below adds the library to a page, and creates a client pointing at the OpenVIVO (<u>http://openvivo.org</u>) TPF endpoint. The client then requests the triples for a particular entity. The triples returned by the query are then used in a second query (.Link) to return all the rdfs:label values from the first query. The .Single function provides a callback function that executes on one returned value when the queries are completed. In the example, the callback displays a single label in the browser console log.

<script href="tpf.js"></script> <script>

const rdfs = "http://www.w3.org/2000/01/rdf-schema#"

Ontology

A small ontology was created to represent the data for the metabolomics consortium application. The ontology simplifies the VIVO ontology and most assertions can be inferred from the VIVO ontology.

For example, in the VIVO ontology, is a person p is the pi of a grant g, the VIVO ontology would assert:

p bearerOf r r a PIrole r relatedBy g

A new object property is PIOF ("is principal investigator of") is used to simplify the above assertions to the equivalent assertion below:

p isPIOf g

A TPF query of the form

p isPIOf *

returns all the grants for which p is the principal investigator. Other "shortcut" object properties include isAuthorOf for publications and isCreatorOf for datasets. Such shortcuts speed up the retrieval of metadata using TPF. The ontology is available here: https://github.com/ctsit/metabolomics_ontology

Progress and Next Steps

TPF is being used to develop web pages for metadata regarding metabolomics coming from multiple sources – Metabolomics Workbench, PubMed, and a spreadsheet describing metabolomics software tools.

The approach described here is easily extended to support:

- Internationalization. The client can accept a language parameter that is used to filter triples to those containing the preferred language, and any secondary languages desired.
- 2. Cross-site TPF. A web page can have any number of clients, each client associated with a specified TPF endpoint. References in one VIVO to triples in a second VIVO can result in the the creation of a second client for the second VIVO and retrieval of desired triples from the

this query returns 3.8M triples. The query

<uri> * *

returns all triples with the specified uri as a subject. When executed on a person's URI in VIVO, the resulting set of triples is the first order graph of the person's information in VIVO.

A TPF endpoint is included as part of the VIVO installation in version 1.10 and above based on the the TPF server available here: https://github.com/LinkedDataFragments

const endpoint = "http://openvivo.org/tpf/core"
const client = new tpf.Client(endpoint)

second VIVO using TPF.

client

.Entity("http://openvivo.org/a/orcid0000-0002-1304-8447")
.Link(rdfs, "label")
.Single(function (label) { console.log(label) })

</script>

Additional functions in the library include: 1) . Results provides a callback that can process a list of returned results. 2) . Query can be used to specify an arbitrary subject, predicate, object TPF query, returning a set of triples. 3) . Type is used to reduce a returned set of triples to entities of a specified rdf:type.

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